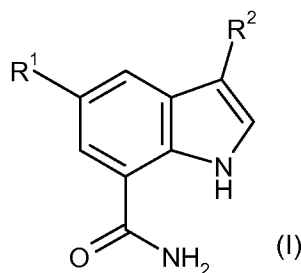


**Amendments to the Claims:**

This listing of claims will replace all prior versions, and listings, of claims in the application:

1. (Previously presented): A compound of Formula (I):



wherein  $R^1$  represents H, halogen, or a group  $-YZ$ ;

Y represents a bond (i.e. is absent),  $C_{1-6}$  alkylene or  $C_{2-6}$  alkenylene;

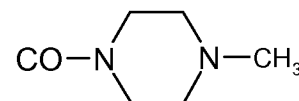
Z represents an aryl or heteroaryl group each comprising 5-14 ring members, said aryl or heteroaryl being optionally substituted by one or more substituents independently selected from halogen, OH,  $C_{1-6}$  alkyl,  $C_{1-6}$  haloalkyl,  $C_{1-6}$  alkoxy,  $C_{1-6}$  haloalkoxy, CN,  $C_{1-6}$  hydroxyalkyl, phenyl,  $O-(CH_2)_{1-6}$ -phenyl,  $NHSO_2R^3$ ,  $NHCOR^3$ ,  $CONR^4R^5$ ,  $SO_2NR^4R^5$ ;

$R^3$ ,  $R^4$  and  $R^5$  independently represent H or  $C_{1-6}$  alkyl;

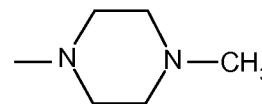
$R^2$  represents a group  $-Y^1Z^1$ ;

$Y^1$  represents a bond (i.e. is absent),  $C_{1-6}$  alkylene,  $C_{2-6}$  alkenylene;

$Z^1$  represents a 6 membered heterocycle which is 4-piperidyl which may be optionally substituted by



one or more substituents independently selected from  $SO_2R^6$ ,  $NHSO_2R^6$ ,  $COR^7$ ,  $NR^7R^8$ ,  $SO_2NR^7R^8$ ,  $C_{1-6}$  alkyl,  $C_{1-6}$  haloalkyl,  $C_{1-6}$  alkoxy,  $C_{1-6}$  haloalkoxy, halogen,  $CONR^7R^8$ ,  $NHCOR^7$ , or phenyl (directly attached or attached by a  $C_{1-6}$ alkylene, CONH,  $C_{2-6}$  alkenylene spacer



and optionally-substituted by one or more substituent selected from  $C_{1-6}$  alkyl,  $C_{1-6}$  haloalkyl,  $C_{1-6}$  haloalkoxy, OH, halogen);

$R^6$  represents H,  $C_{1-6}$  alkyl,  $-(CH_2)_n$  phenyl or  $-(CH_2)_n$  naphthyl (where n is 0 or 1 and each of which phenyl or naphthyl may be optionally substituted by one or more substituents independently selected from  $C_{1-6}$  alkyl,  $C_{1-6}$  alkoxy, halogen,  $NR^7R^8$ ,  $C_{1-6}$  haloalkyl,  $C_{1-6}$  haloalkoxy), CN or  $-(O)_p$  phenyl (where p is 0 or 1 and the phenyl is optionally substituted by one or more substituents independently selected from halogen,  $C_{1-6}$  alkyl or  $C_{1-6}$  alkoxy));

$R^7$  and  $R^8$  independently represents  $C_{1-6}$  alkyl, H,  $C_{1-6}$  alkylene  $NR^9R^{10}$ ; and

$R^9$  and  $R^{10}$  independently represents  $C_{1-6}$  alkyl, H;

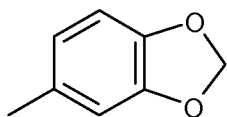
or a pharmaceutically acceptable salt thereof.

2. (Original): A compound according to claim 1 wherein  $R^1$  is YZ.

3. (Original): A compound according to claim 2 wherein Y is a bond or  $-CH = CH-$ .

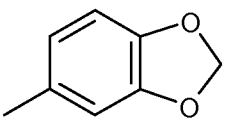
4. (Original): A compound according to claim 3 wherein Y is a bond.

5. (Previously presented): A compound according to claim 1 wherein Z is phenyl (which may be unsubstituted or substituted once or twice by substituents independently selected from  $C_{1-3}$  alkoxy, CN, OH, phenyl,  $-OCH_2$  phenyl  $NHSO_2R^3$ ,  $NHCO_2R^3$ ,  $CONR^4R^5$ ,  $SO_2NR^4R^5$ , halogen,  $C_{1-3}$  hydroxyalkyl,  $C_{1-4}$  alkyl) or a heteroaryl group selected from benzofuranyl, quinolinyl,



, pyrimidinyl, thiophenyl, isoxazolyl, pyridinyl (each of which may be optionally substituted by one or two groups independently selected from  $C_{1-3}$  alkyl,  $C_{1-3}$  alkoxy, halogen).

6. (Original): A compound according to claim 5 wherein Z is phenyl (which is unsubstituted or substituted once by a substituent selected from phenyl,  $OCH_2$  phenyl,  $NHSO_2CH_3$ ,  $NHCOCH_3$ ,  $CONH_2$ ,  $CON(CH_3)_2$ , Cl, F,  $OCH_3$ , CN, OH,  $CH_2OH$ ,  $CH_3$ ,  $C(CH_3)_3$ ) or a heterocyclic group selected

from benzofuranyl, quinolinyl, , pyrimidinyl, thiophenyl, benzothiophenyl, isoxazolyl, pyridinyl (each of which is substituted or is substituted once by a group selected from -OCH<sub>3</sub>, CH<sub>3</sub>, F).

7. (Original): A compound according to claim 6 wherein Z is phenyl (which is unsubstituted or substituted once by a substituent selected from phenyl, OCH<sub>2</sub> phenyl, NHSO<sub>2</sub>CH<sub>3</sub>, NHCOCH<sub>3</sub>, CONH<sub>2</sub>, CON(CH<sub>3</sub>)<sub>2</sub>, Cl, F, OCH<sub>3</sub>, CN, OH, CH<sub>2</sub>OH, CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>).

8. (Original): A compound according to claim 7 wherein Z is phenyl.

Claims 9-10 (Cancelled)

11. (Previously presented): A compound according to claim 1, wherein Y<sup>1</sup> is a bond or C<sub>1-3</sub> alkylene.

Claim 12 (Cancelled)

13. (Previously presented): A compound according to claim 1, wherein Z<sup>1</sup> is a 6 membered heterocycle which is 4-piperidyl substituted by SO<sub>2</sub>R<sup>6</sup>.

Claims 14-18 (Cancelled)

19. (Previously presented): A pharmaceutical composition, comprising a compound as claimed in claim 1, or a pharmaceutically acceptable salt thereof and one or more of pharmaceutically acceptable carriers, diluents and excipients.

Claims 20-41 (Cancelled)